# ANALYTICAL RESULTS OF SURFACE WATER SAMPLES COLLECTED FROM RACCOON CREEK October 20, 1999 Sampling Event

### Prepared for:

LYONDELL CHEMICAL WORLDWIDE, INC./BEAZER EAST INC.

### Prepared by:

Applied Hydrology Associates, Inc. Pittsburgh, Pennsylvania Denver, Colorado

November 10, 1999



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#### 1.0 INTRODUCTION

This report presents the results of surface water samples collected from Raccoon Creek at the Lyondell Chemical Worldwide, Inc. / Beazer East Inc. Monaca, PA site during the October 20, 1999 quarterly monitoring event. The samples were collected in compliance with Appendix D of the 1997 Consent Order and Agreement (1997 CO&A) between ARCO Chemical Company<sup>1</sup>, BEI and the Pennsylvania Department of Environmental Protection (PADEP) dated October 20, 1997.

#### 2.0 SAMPLING

Surface water samples were collected at Transect E as defined in the 1997 CO&A. The location of Transect E is shown in Figure 1. In addition, water elevations were measured in nearby monitoring wells and the results are presented in Appendix A.

A total of eight surface water samples, including a duplicate, were collected from Raccoon Creek on October 20, 1999. These samples were collected at the same three locations along Transect E as in previous sampling events. The locations are shown in Figure 2 and are at the center of the stream, and approximately 30 feet from the east and west banks. At the center location, samples were collected at three depths; 6 inches below the surface, 2 inches above the bottom, and midway between the surface and bottom. Samples from the east and west sides of the transect were collected at two depths; 2 inches above the bottom, and midway between the surface and bottom.

During sampling a boat was stationed at Transect E using a rope secured to the east and west shores of Raccoon Creek. The samples were collected by using a peristaltic pump to pump water from the desired depth into three 40-ml vials preserved with hydrochloric acid. Samples were collected from the required depths utilizing tubing secured to a vertical steel rod lowered from the boat until it rested on the bottom of the creek. The rod did not penetrate the sediments on the creek bottom because a 1-foot diameter disc constructed of steel mesh is fastened perpendicular to the bottom of the rod.

Two tubes were used. The bottom of the "deep sample tube" was secured to the probe 2 inches from the bottom of the probe. The bottom of the "mid-depth sample tube" is adjustable and was secured to the probe mid-depth at each location. Care was taken not to disturb the sediments at the sampling location and the pumped water was closely monitored to ensure sediment was not included in the sample. One gallon of water was pumped through the tubing before each sample was obtained in order to purge the tubing.

<sup>&</sup>lt;sup>1</sup> ARCO Chemical Company is now Lyondell Chemical Worldwide

The samples were uniquely numbered as follows to identify the location, depth and date of sampling:

RC-EC-00-1099

Where:

RC indicates Raccoon Creek

EC indicates Transect E and location (C=Center, L = Left bank, R = Right bank

[facing downstream])

00 indicates sample depth in feet and tenths of a foot (0.0 feet)

indicates the month and year collected (October 1999)

Samples were logged onto a chain of custody form (included in of the Analytical Report in Appendix B) and stored on ice until receipt by Reliance Laboratories Inc. in Edison, NJ. Reliance analyzed the samples using USEPA Method 524.2 for BTEXS.

#### 3.0 RESULTS

The analytical results are presented in Table 1. Benzene was detected in six of the seven locations and concentrations in samples where benzene was detected ranged from 0.18  $\mu$ g/L in Sample RC-EL-00-1099<sup>2</sup> to 0.43  $\mu$ g/L in sample RC-EC-00-1099. Sampling locations and depths are shown on Figure 2, along with the concentration of benzene at each location. Water levels in wells near Raccoon Creek are presented in Appendix A.

Table 1
Summary of Analytical Results for Samples Collected from Raccoon Creek

Sample Name	Benzene	Toluene	Ethylbenzene	Xylene	Styrene
RC-EL-00-1099	0.14	0.60	<0.22	0.60	< 0.58
RC-EL-00-1099A	0.21	1.26	0.25	1.16	< 0.58
RC-EL-19-1099	0.27	1.32	0.28	1.4	< 0.58
RC-EC-00-1099	0.43	2.49	0.48	3.11	< 0.58
RC-EC-33-1099	0.42	2.18	0.43	2.54	< 0.58
RC-EC-69-1099	0.20	0.6	<0.22	0.67	< 0.58
RC-ER-00-1099	0.23	1.03	<0.22	1.05	< 0.58
RC-ER-65-1099	< 0.13	<0.6	<0.22	0.47	< 0.58

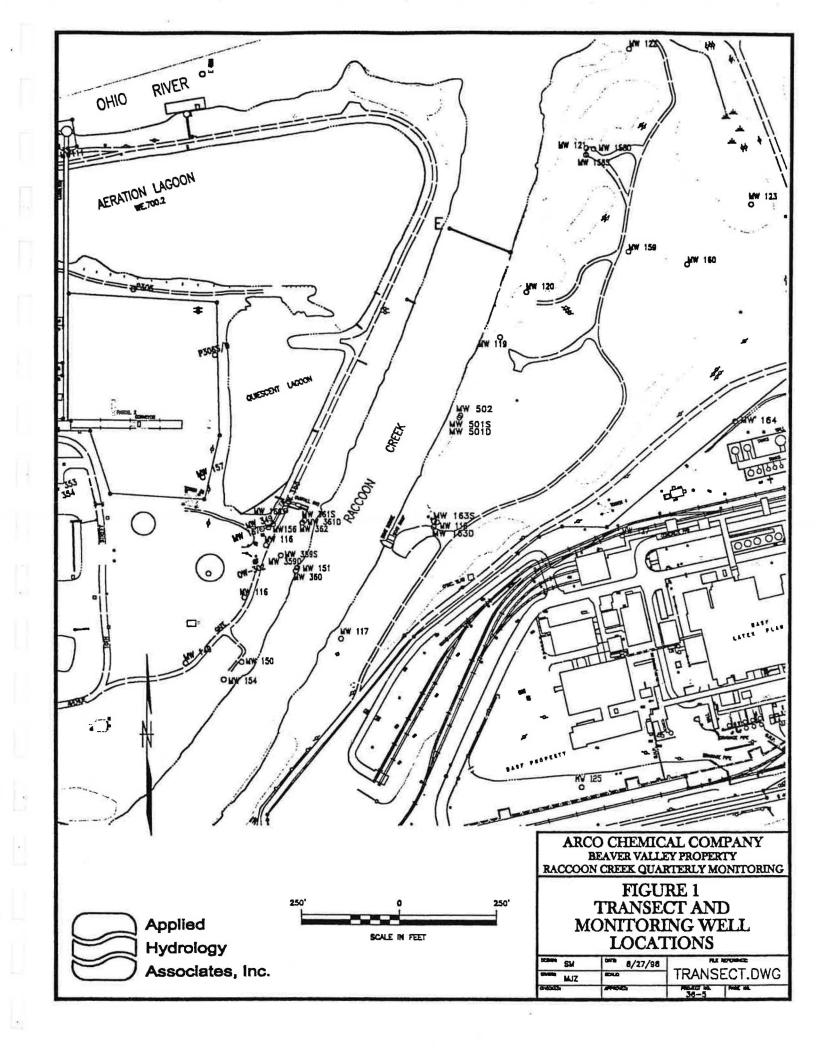
The analytical data were validated upon receipt and found to be acceptable. A Data Validation Report is provided in Appendix B. Table 2 presents the historical concentration of benzene in Raccoon Creek at Transect E during all monitoring events to date.

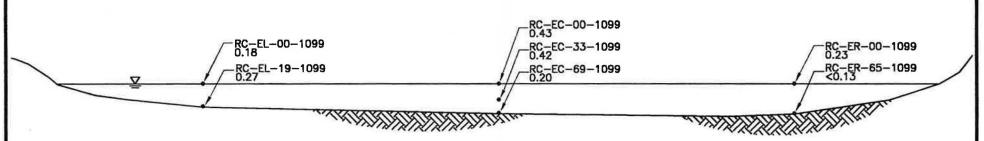
<sup>&</sup>lt;sup>2</sup> The benzene concentration of 0.18 ug/L is the average of the concentration in the two duplicates samples, RC-EL-00-1099 and RC-EL-00-1099A.

### Historic Benzene Concentrations at Transect E (ug/L)

Sampling Location	Sampling Depth	7/23/97	10/28/97	2/25/98	5/21/98	7/29/98	10/27/98	2/3/99	4/27/99	7/22/99	10/20/99
30 Feet off West Bank	Mid-depth	0.28	<0.13	<0.13	0.70	<0.13	1.57 <sup>(1)</sup>	0.37	< 0.66	<0.13	0.18(1)
30 Feet off West Bank	Deep	0.81	<0.13	<0.13	0.70	<0.13	0.61 <sup>(1)</sup>	0.49	< 0.66	<0.13	0.27
Center of Creek	Shallow	0.24	<0.13	0.38	0.70	<0.13	<0.13	0.61(1)	< 0.66 <sup>(1)</sup>	<0.13 <sup>(0)</sup>	0.43
Center of Creek	Mid- Depth	0.18	<0.13	0.49	0.64	<0.13	0.2	0.64	< 0.66	<0.13	0.42
Center of Creek	Deep	0.46	<0.13	0.30	0.60	<0.13	<0.13	0.69	< 0.66	<0.13	0.20
30 Feet off East Bank	Mid-depth	0.16	<0.13	<0.13	<0.13	0.13	0.52	< 0.13	< 0.66	<0.13	0.23
30 Feet off East Bank	Deep	<0.13	<0.13	0.14	0.22	0.22	<0.13	< 0.13	< 0.66	<0.13	<0.13

<sup>(1)</sup> Results shown are the average of the blind duplicate samples.





## CREEK SECTION LOOKING DOWNSTREAM

### **LEGEND**

• SURFACE WATER SAMPLE LOCATION
ALL CONCENTRATIONS IN ug/L



Applied
Hydrology
Associates, Inc.



LYONDELL CHEMICAL WORLDWIDE BEAVER VALLEY PROPERTY RACCOON CREEK QUARTERLY MONITORING

FIGURE 2

SURFACE WATER
BENZENE CONCENTRATIONS
AT TRANSECT "E"

OCTOBER 20, 1999

SM	B/17/98	PLE REPORTED				
JLS	NOT TO SCALE	BENZENE.dwg				
Odoba	APPROPRIE	38-5 Polit Hd.				

### Appendix A

Groundwater Elevations, East and West Sides of Raccoon Creek

## GROUNDWATER LEVELS ON THE EAST AND WEST SIDES OF RACCOON CREEK OCTOBER 20, 1999

Well Number			Depth to Water	Table 1	Calculated SPL	Comments
	(TOC)	from TOC (2)	from TOC (2)	Level Elevation (1)	Thickness (3)	
	(ft. amsl)	(ft. amsl)	(ft. amsl)	(ft. amsl)	(ft. amsl)	
		N	Ionitoring V	Vells Screened	in Silty Clay	Unit
			7.00	OTH AREA		
MW - 360	685.84	ND	3.12	682.72	N/A	
MW - 170	706.70	ND	22.44	684.26	N/A	
MW - 362	689.43	ND	6.29	683.14	N/A	
			F	RACCOON CREEK	AREA	
						Monitoring well is dry. Screened above water table
MW- 118	690.39	ND	7.06	683,33	N/A	Bottem of well is 7.06 feet below TOC.
MW - 502	701.86	ND	19.08	682.78	N/A	
MW - 119	705.59	ND	22.77	682.82	N/A	
MW - 120	709.42	ND	26.50	682.92	N/A	
MW - 121	713.90	ND	30.91	682.99	N/A	
MW - 152	696.35	ND	13.80	682.55	N/A	
		Monitor	ing Wells So	reened in Upp	er Sand and	Gravel Unit
	****	112011101	ing Weils St	OTH AREA	or Sunt und	GIATOL CHIV
MW - 344	709.42	ND	24.89	684.53	N/A	
MW - 359S	692.93	ND	9.50	683.43	N/A	
MW - 361S	689.40	ND	6.23	683.17	N/A	
						Well had excessive pressure build up from OTH
MW - 169	707.93	ND	25.35	682.58	N/A	Sprage event (7/13/99 - 7/14/99).
						Top of casing changed from 707.36 to 711.06 on 11/98
						accommodate respiration monitoring well head.
MW - 167	711.06	ND	28.00	683.06	N/A	Monitoring well stick up is 3.70 above orig. TOC
7 CO C	<b>600.05</b>			RACCOON CREEK		
MW - 163S	690.87	ND	7.81	683.06	N/A	
MW - 501S	701.30	ND	18.72	682.58	N/A	
MW - 162S	706.05	ND	23.22	682.83	N/A	
MW - 159	708.99	ND	25.99	683.00	N/A	
MW - 160	701.00	ND	18.01	682.99	N/A	
MW - 158S	713.60	ND	30.62	682.98	N/A	
MW - 122	692.78	ND	9.81	682.97	N/A	
Note: See figur		T1 4' 6T	00 : 7 1	TILL C. TOO		
				to Water from TOC.	1	
				e. ND means no SPL		s not applicable, no SPL was detected.

## GROUNDWATER LEVELS ON THE EAST AND WEST SIDES OF RACCOON CREEK OCTOBER 20, 1999

Well Number	Top of Casing (TOC) (ft. amsl)	Depth to SPL from TOC (2) (ft. amsl)	Depth to Water from TOC (2) (ft. amsl)	Calculated Water Level Elevation (1) (ft. amsl)	Calculated SPL Thickness (3) (ft. amsl)	Comments
		Monitori	ng Wells Sc	reened in Lowe	er Sand and (	Gravel Unit
MW 345	708.91	ND	25.84	683.07	N/A	
MW 361D	689,35	ND	6.30	683.05	N/A	
MW 359D	692.80	ND	9.70	683.10	N/A	
MIV JUST	072.00	1,12		RACCOON CREEK		
MW 163D	689.62	ND	6.54	683.08	N/A	
MW 501D	701.44	ND	18.50	682.94	N/A	
MW 166D	703.95	ND	20.98	682.97	N/A	
MW 158D	712.04	ND	29.16	682.88	N/A	
		W	ater Levels i	n Raccoon Cre	ek and Ohio	River
			RACCOO	N CREEK AREA S	TAFF GAUGE	
Time of	Staff Gauge Elevation (ft. amsl)	Staff Gauge	Calculated Water Level Elevation			Comments
Observation	(4) (5)	Reading	(ft. amsl)			
10:33	685.00	0.70	682.70			
13:00	685.00	0.70	682.70			ř.
				IO RIVER. STAFF	GAUGE	
9:46	685.96	2.75	682.71			*
12:30	685.96	2.75	682.71			
						×
Note: See figure		ElAi - CT/	20 minus 20 mil	TOC		-9-
				to Water from TOC.	rupp detected	
				e. ND means no SPL		not applicable, no SPL was detected.
					IOC. N/A means	not applicable, no SPL was detected.
4) Elevation 68	5 00 is agricult	ant to 3 AA mode	on chaff gamas at	Daccoon Crook		

# Appendix B Data Validation Report



1200 South Parker Road, Suite 100

Denver, CO 80231

Tel: (303) 873-0164

Fax: (303) 873-6110

### MEMORANDUM

TO:

Files

FROM:

Skip Meier, Applied Hydrology Associates

DATE:

November 11, 1999

SUBJECT:

Data Validation Results, Lyondell Chemical Worldwide Beaver Valley Property

Data validation was performed on the volatile organic analytical data from eight surface water samples obtained from Raccoon Creek on October 20, 1999 and also on a Rinsate Blank and Trip Blank. The validation was performed in accordance with the "USEPA Contract Laboratory Program National Functional Guidelines for Organic Data Review". Reliance Laboratories Inc. performed the analysis using EPA Method 524.2. The samples reviewed included:

Field Sample ID	Lab Sample ID
RC-EL-00-1099	R-6204.4
RC-EL-00-1099A	R-6204.8
RC-EL-19-1099	R-6204.3
RC-EC-00-1099	R-6204.7
RC-EC-33-1099	R-6204.6
RC-EC-69-1099	R-6204.5
RC-ER-00-1099	R-6204.2
RC-ER-65-1099	R-6204.1
Rinsate Blank	R-6204.9
Trip Blank	R-6204.10

Items reviewed and actions taken were as follows:

#### √ Method:

The ten samples were analyzed for BTEXS by method USEPA 524.2 on October 21, 1999.

### Holding Time:

All Samples were analyzed within the 14-day holding time.

#### √ <u>Blanks:</u>

No target compounds were detected in the associated method blank.

#### **√** System Monitoring Compounds:

The "USEPA Contract Laboratory Program National Functional Guidelines for Organic Data Review" indicate that "Recoveries for system monitoring compounds in volatile samples and blanks must be within the limits specified in the Method." However, Method 524.2 does not specify a required recovery. Nevertheless, 4-

bromofluorobenzene and 1,2-dichlorobenzene-d4 surrogate recoveries were within 93-108 percent and this is acceptable.

√ Internal Standards:

All fluorobenzene internal standards were within the established criteria for area internal standard and retention time.

√ GC/MS Instrument Performance Check:

All bromofluorobenzene (BFB) tunes met the ion abundance criteria. Analysis of the instrument performance check solution was performed at the beginning of each 12-hr period during which the samples were analyzed.

√ <u>Initial Calibrations:</u>

The initial calibration performed on October 21, 1999 for Instrument HP5971A met the 30 percent relative standard deviation (RSD) and 0.05 minimum relative response factor criteria for all compounds.

√ Continuing Calibrations:

Continuing calibration was run and compared to the correct initial calibration. All continuing calibrations met the 25 percent difference and minimum relative response factor criteria for all compounds.

√ Matrix Spike/Duplicate:

The matrix spike/duplicate results for recovery and RPD were within the Quality Control limits.

### √ Target Compound Indentification/Quantitation:

No problems were identified with compound identification or quantities.

√ Field Duplicate:

A field duplicate was collected during this sampling event. The duplicate sample was denoted by an "A" at the end of the sample name. The pair is RC-EL-00-1099 and duplicate RC-EL-00-1099A. Table 1 below summarizes the RPD for the sample/duplicate pair<sup>1</sup>.

Table 1: Relative Percent Difference (RPD)

Sample Name	Benzene (ppb)	RPD (%)	Toluene (ppb)	RPD (%)	Ethyl- Benzene (ppb)	RPD (%)	Xylene (ppb)	RPD (%)	Styrene (ppb)	RPD (%)
RC-EL-00-1099 RC-EL-00-1099A	14 21	40	0.60 1.26	71	<0.22 0.25	NA	0.60 1.16	64	ND ND	NA NA

ND = Non Detect

NA = Not Applicable

√ Summary:

No inconsistencies were noted except that poor agreement was seen between the duplicate sample pair RC-EL-00-1099 and RC-EL-00-1099A. (See Table 1). No BTEXS compounds were detected in either the trip blank or the field blank.

The equation for calculating RPD is:  $_{R}$   $_{P}$   $_{D}$  = 2 \*  $\frac{|S - D|}{|S + D|}$  \* 1 0 0 where S = sample concentration and D

<sup>=</sup> duplicate concentration

### RELIANCE LABORATORIES, INC.



175 MAY STREET, EDISON, NJ 08837 PH (732) 738-5454 FAX (732) 738-5841 EMAIL: 74201.3501@COMPUSERVE.COM

### **ANALYTICAL REPORT**

For Lyondell Chemical Monaca, PA 15061

Project: Raccoon Creek

### RELIANCE LABORATORIES, INC.



175 MAY STREET, EDISON, NJ 08837 PH (732) 738-5454 FAX (732) 738-5841 EMAIL: 74201.3501@COMPUSERVE.COM

### **ANALYTICAL DATA REPORT**

for

Lyondell Chemical Monaca, PA 15061 Project: Raccoon Creek

Date Received: 10/21/99

Sample ID	Lab ID #
:	
RC-ER-65-1099	R-6204.1
RC-EF-00-1099	R-6204.2
RC-EL-19-1099	R-6204.3
RC-EL-00-1099	R-6204.4
RC-EC-69-1099	R-6204.5
RC-EC-33-1099	R-6204.6
RC-EC-00-1099	R-6204.7
RC-EL-00-1099A	R-6204.8
Rinsate Blank	R-6204.9
Trip Blank	R-6204.10

These samples have been analyzed by EPA Method 524.2 for a selected compound list. The results are not designed for use for drinking water purposes.

G. P. Kirpalani

Manager

GPK/vb

### RELIANCE LABORATORIES INC.



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### RELIANCE LABORATORIES, INC.



175 MAY STREET, EDISON, NJ 08837 PH (732) 738-5454 FAX (732) 738-5841 EMAIL: 74201.3501@COMPUSERVE.COM

### LABORATORY CHRONICLE

Date Received: 10/21/99 Date Sampled: 10/20/99 Sample ID: As per chain of custody	
Organic Extraction:	
1 Acids	
2 Base / Neutrals 3 Pesticides/PCB's 4 TPHC	
Analysis:	
1 Volatiles 10/21/99	
2 Acids	
3 Base/Neutrals	
4 Pesticides/PCB's	
5 TPHC	
Inorganics:	
1 Metals	
2 Cyanides	
3 Phenols	
Other Analysis:	
Supervisor Review & Approval	

RELIANCE LABORATORIES INC.



3090 WOODBRIDGE AVENUE, EDISON NJ 08837 PH (908) 738-5454 FAX (908) 738-5841

### **NON-CONFORMANCE SUMMARY**

Reliance Labs received 8 water sample, 1 Rinsate Blank and 1 Trip blank for BTEXS by method (EPA 524.2) from Lyondell Chemical on 21 October 1999. Samples consisted of 10 vials.

Matrix spike recovery analysis was performed on samples and results are attached. 10mL of sample was used on spike/spike dup since not enough sample was provided

All analyses were performed within the required holding time.

### 4. Run Sequence

- A. Tune Instrument
- 1. Inject 1µL of 25 ppm BFB into GC.
  - a. Tune must pass against criteria.
  - b. Tune must be run before any samples, blank or calibrations can be run.
  - c. From time to tune 12 hours are available to run all QC data and samples.

### B. Five Point Calibration Curve

- 1. Purge five (5) concentrations of standard solutions containing all the target analysis at 1 ppb, 2 ppb, 5 ppb, 10 ppb, and 20 ppb.
- 2. The above standard must be run within 12 hours of injecting the BFB tune.
- 3. Created a calibration curve with the above standard runs.
  - a. If the 30% RSD deviation is exceeded the standards must be run again (still within 12 hours)
- 4. Create an identification file from this calibration curve for automated quantification.
- C. If time remains in the 12-hour run period go to step F.
- D. If the 12-hour period has expired, a new tune must be injected and a new sequence must be started.
- E. Once an initial calibration curve is established a continuing calibrations check may be run. A continuing calibration check is required every time the mass spectrometer is tuned.
  - 2 ppb concentration of all target compounds is purged and quanted against current ID file.
  - Check the response factors of this run against the average RF of the calibration file. The RF of the continuing calibration must be within ± 50% D (difference) of the 5 point for all compounds.
  - 3. The area counts of internal standard and surrogates must not be decreased by >30% from the most recent continuing calibration standard nor decrease by >50% from the initial calibration standard.

### F. Daily Blank

- 1. Purge 20 ml of laboratory reagent water (nanopure) with 5 ppb internal standard and 5 ppb each surrogate.
- 2. Run this blank and quant against current ID file.
- 3. If blank does not meet criteria, it must be rerun before analyzing any samples.

### G. Samples

- 1. Fill 25 ml syringe until it overflows with sample. Then adjust the volume to 20 ml exactly.
- 2. Inject 5 µl each 25 ppm internal standard and surrogate standard solution into each sample.
- Run and quant against the current 5 point calibration curves.
- 4. Any sample with target compound over 50 ppb must be rerun at the appropriate dilution.
- 5. Any sample not injected in 12-hour period must be rerun.

### H. Quality Control Sample (QCS)

1. Analyze a QCS from an external source at least quarterly.

### STANDARD OPERATING PROCEDURE METHOD 524.2

1. Scope
This is the general method for the procedure used to identify purgeable volatile organics in portable water. The sample is purged with ultra high purity helium and concentrated into a trap. The volatiles are then thermally desorbed onto a megabore column and identified using a mass

- 2. Equipment and Apparatus
- A. Sample containers- 40ml screw caps amber vials.
- B. Purge and Trap System.
  - 1. 25cm VOCARB 3000 trap.
- C. Glassware

spectrometer detector.

- 1. 20 ml fritted purging vessels.
- 2. 25 ml teflon sealed syringe with lever lock assembly.
- 3. 10 μL syringes.
- D. Gas Chromographic / Mass Spectrometer.
  - 1. Column type J&W

75 m, 0.53 mm ID, DB624 3 microns

- E. Apparatus Conditions
  - 1. Tekmar (purge and trap)

a. Purge time : 2 min.

b. Desorb time and temp.: 250° for 2 min.

c. Bake time and temp. : 260° for 12 min.

d. Flow rate : 15 cc/min.

2. GC Conditions

a. Column flow 15 cc/min.

b. Initial temp. 35° C

c. Ramping Rate 6° C/min. d. Final temp. 200° C

e. Run time 47.25 min.

Initial time 6 min.

- 3. Stock Standards
- A. Internal Standard
  - 1. Flourobenzene

f.

- B. Surrogates
  - 1. 1.2-dichlorobenzene-d4
  - 2. 4-bromoflurobenzene
- C. Prepare standard solutions for all target compounds and surrogates at 20 ppm.
- D. Prepare internal standard at 20 ppm in methanol.
  - 1. Prepare all standards and store in teflon sealed 1 ml vials.

### R E L I A N C E LABORATORIES INC.



175 MAY STREET, EDISON, NJ 08837 PH (908) 738-5454 FAX (908) 738-5841 EMAIL: 74201.3501@COMPUSERVE.COM

LABORATORY ID NJ DEP NO. 12687 PA DER NO. 68437

#### **CERTIFICATE OF ANALYSIS**

Customer:

Lyondell Chemical

Sample:

**Aqueous Samples** 

Date Sampled:

20 October 1999

Lab ID:

R-6204

Reference:

AHA / Monaca

22 October 1999

Units: μg/L

Sample ID	Benzene	Toluene	Ethylbenzene	Xylene	Styrene
RC-ER-65-1099 RC-ER-00-1099 RC-EL-19-1099 RC-EL-00-1099 RC-EC-69-1099 RC-EC-33-1099 RC-EC-00-1099	< 0.13 0.23 0.27 0.14 0.20 0.42 0.43	< 0.6 1.03 1.32 0.60 0.60 2.18 2.49	< 0.22 < 0.22 0.28 < 0.22 < 0.22 0.43 0.48	0.47 1.25 1.40 0.60 0.67 2.54 3.11 1.16	< 0.58 < 0.58 < 0.58 < 0.58 < 0.58 < 0.58 < 0.58 < 0.58
RC-EL-00-1099A Rinsate Blank Trip Blank	0.21 < 0.13 < 0.13	1.26 < 0.6 < 0.6	0.25 < 0.22 < 0.22	< 0.22 < 0.22	< 0.58 < 0.58

G. P. Kirpalani Manager

Data File : C:\HPCHEM\1\DATA\V6633.D Acq On : 21 Oct 99 7:46 pm

Vial: 3 Operator: vb

Inst : 5971 - In

Sample : R-6204.1 Misc : Lyondell - RC-ER-65-1099

Multiplr: 1.00

Quant Time: Oct 22 11:27 1999

Method

: C:\HPCHEM\1\METHODS\RUN524.M

Title

: 524.2 Purgable Organics

Last Update : Thu Oct 21 13:22:11 1999 Response via : Multiple Level Calibration

Internal Standards	R.T.	QIon	Response	Conc Units	Dev(Min)
1) Fluorobenzene	12.49	96	506914	5.00 ug/L	-0.04
System Monitoring Compounds 43) 4-bromofluorobenzene 55) 1,2-dichlorobenzene-d4	25.65 30.85	95 152	191281 126071	%R 5.02 ug/L 5.13 ug/L	100.39% 102.69%
Target Compounds 26) Toluene 36) m&p-xylenes 37) o-xylene	17.39 22.62 23.91	91 106 91	59835 25484 31975	0.38 ug/L 0.21 ug/L 0.26 ug/L	Qvalue 92 74 93

<sup>(#) =</sup> qualifier out of range (m) = manual integration V6633.D RUN524.M Fri Oct 22 11:27:38 1999

Vial: 3

Multiplr: 1.00

: 5971 - In

Operator: vb

Inst

Data File : C:\HPCHEM\1\DATA\V6633.D

Acq On : 21 Oct 99 7:46 pm

: R-6204.1 Sample

Method

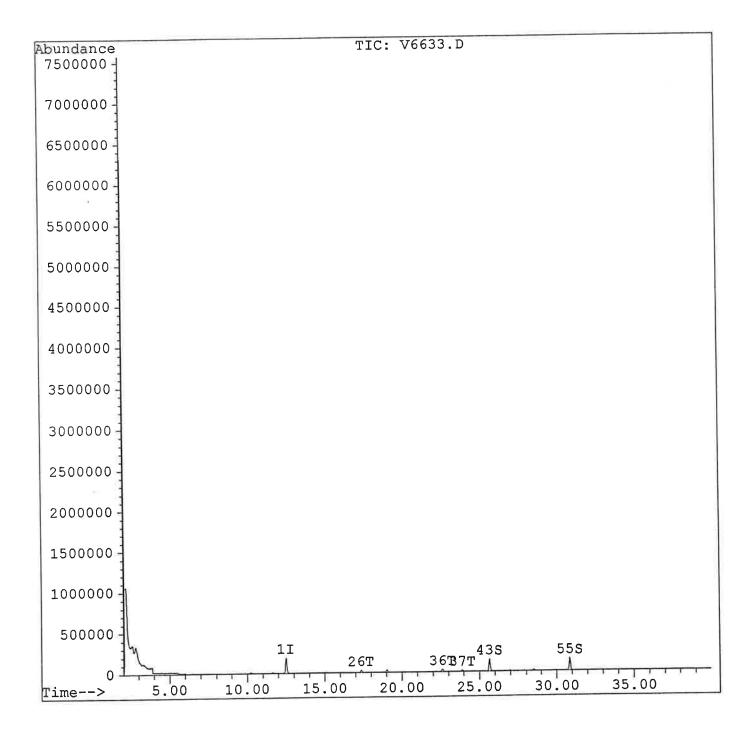
Title

Misc

Quant Time: Oct 22 11:27 1999

: Lyondell - RC-ER-65-1099

: C:\HPCHEM\1\METHODS\RUN524.M : 524.2 Purgable Organics Last Update : Thu Oct 21 13:22:11 1999 Response via : Multiple Level Calibration



Data File : C:\HPCHEM\1\DATA\V6627.D Acq On : 21 Oct 99 3:08 pm

Vial: 11 Operator: vb

Inst : 5971 - In

Multiplr: 1.00

Sample : R-6204.2 Misc : Lyondell - RC-ER-00-1099

Quant Time: Oct 22 11:38 1999

Method : C:\HPCHEM\1\METHODS\RUN524.M
Title : 524.2 Purgable Organics Last Update : Thu Oct 21 13:22:11 1999 Response via : Multiple Level Calibration

Internal Standards	R.T.	QIon	Response	Conc Units	Dev(Min)
1) Fluorobenzene	12.52	96	487047	5.00 ug/L	-0.01
System Monitoring Compounds 43) 4-bromofluorobenzene 55) 1,2-dichlorobenzene-d4	25.67 30.88	95 152	188380 127564	%I 5.14 ug/L 5.41 ug/L	Recovery 102.90% 108.14%
Target Compounds 19) Benzene 26) Toluene 35) Ethylbenzene 36) m&p-xylenes 37) o-xylene	11.68 17.41 22.22 22.65 23.91	78 91 91 106 91	41247 156995 35940 65783 82027	0.23 ug/L 1.03 ug/L 0.21 ug/L 0.56 ug/L 0.69 ug/L	Qvalue 93 99 95 96 94

<sup>(#) =</sup> qualifier out of range (m) = manual integration V6627.D RUN524.M Fri Oct 22 11:38:14 1999

Vial: 11 Operator: vb

Multiplr: 1.00

: 5971 - In

Inst

Data File : C:\HPCHEM\1\DATA\V6627.D

Acq On : 21 Oct 99 3:08 pm

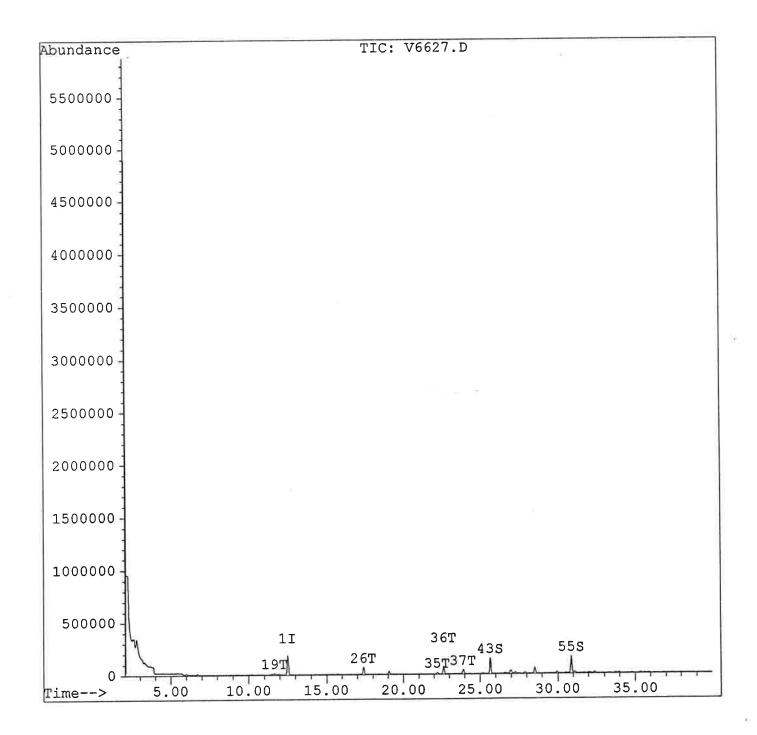
Sample : R-6204.2

Misc : Lyondell - RC-ER-00-1099

Quant Time: Oct 22 11:38 1999

Method : C:\HPCHEM\1\METHODS\RUN524.M

Title : 524.2 Purgable Organics
Last Update : Thu Oct 21 13:22:11 1999
Response via : Multiple Level Calibration



Data File : C:\HPCHEM\1\DATA\V6628.D Vial: 12 Acq On : 21 Oct 99 3:54 pm Operator: vb

Inst : 5971 - In

Sample : R-6204.3 Misc : Lyondell - RC-EL-19-1099 Multiplr: 1.00

Quant Time: Oct 22 11:36 1999

Method : C:\HPCHEM\1\METHODS\RUN524.M
Title : 524.2 Purgable Organics Last Update : Thu Oct 21 13:22:11 1999 Response via : Multiple Level Calibration

Internal Standards	R.T.	QIon	Response	Conc Units	Dev(Min)
1) Fluorobenzene	12.52	96	473211	5.00 ug/L	0.00
System Monitoring Compounds 43) 4-bromofluorobenzene 55) 1,2-dichlorobenzene-d4	25.68 30.88	95 152	173496 114035	%! 4.88 ug/L 4.97 ug/L	Recovery 97.54% 99.50%
Target Compounds 19) Benzene 26) Toluene 35) Ethylbenzene 36) m&p-xylenes 37) o-xylene	11.67 17.40 22.22 22.64 23.92	78 91 91 106 91	47696 196163 46104 70813 89387	0.27 ug/L 1.32 ug/L 0.28 ug/L 0.62 ug/L 0.78 ug/L	Qvalue 90 97 98 99

<sup>(#) =</sup> qualifier out of range (m) = manual integration V6628.D RUN524.M Fri Oct 22 11:36:48 1999

Vial: 12

: 5971 - In

Operator: vb

Multiplr: 1.00

Inst

Data File : C:\HPCHEM\1\DATA\V6628.D

: 21 Oct 99 Acq On

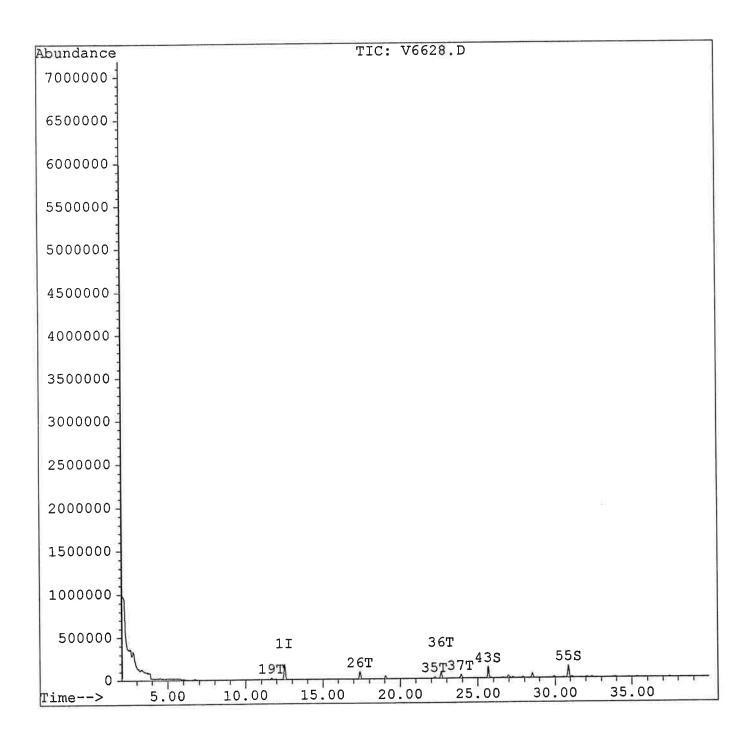
: R-6204.3 Sample

: Lyondell - RC-EL-19-1099 Misc

Quant Time: Oct 22 11:36 1999

: C:\HPCHEM\1\METHODS\RUN524.M

Method : 524.2 Purgable Organics Title Last Update : Thu Oct 21 13:22:11 1999 Response via : Multiple Level Calibration



Data File: C:\HPCHEM\1\DATA\V6629.D Vial: 13
Acq On: 21 Oct 99 4:41 pm Operator: vb

Sample : R-6204.4 Inst : 5971 - In

Misc : Lyondell - RC-EL-00-1099 Multiplr: 1.00

Quant Time: Oct 22 11:35 1999

Method : C:\HPCHEM\1\METHODS\RUN524.M

Title : 524.2 Purgable Organics
Last Update : Thu Oct 21 13:22:11 1999
Response via : Multiple Level Calibration

Internal Standards	R.T.	QIon	Response	Conc Units	Dev(Min)
1) Fluorobenzene	12.52	96	524441	5.00 ug/L	0.00
System Monitoring Compounds 43) 4-bromofluorobenzene 55) 1,2-dichlorobenzene-d4	25.66 30.88	95 152	198829 133554	%! 5.04 ug/L 5.26 ug/L	Recovery 100.86% 105.15%
Target Compounds 19) Benzene 26) Toluene 35) Ethylbenzene 36) m&p-xylenes 37) o-xylene	11.68 17.40 22.25 22.64 23.93	78 91 91 106 91	27151 99495 20370 30845 46195	0.14 ug/L 0.60 ug/L 0.11 ug/L 0.24 ug/L 0.36 ug/L	Qvalue 90 98 94 96

<sup>(#) =</sup> qualifier out of range (m) = manual integration V6629.D RUN524.M Fri Oct 22 11:35:13 1999

Data File : C:\HPCHEM\1\DATA\V6629.D

Acq On : 21 Oct 99 4:41 pm

Vial: 13
Operator: vb

Sample : R-6204.4

Inst : 5971 - In

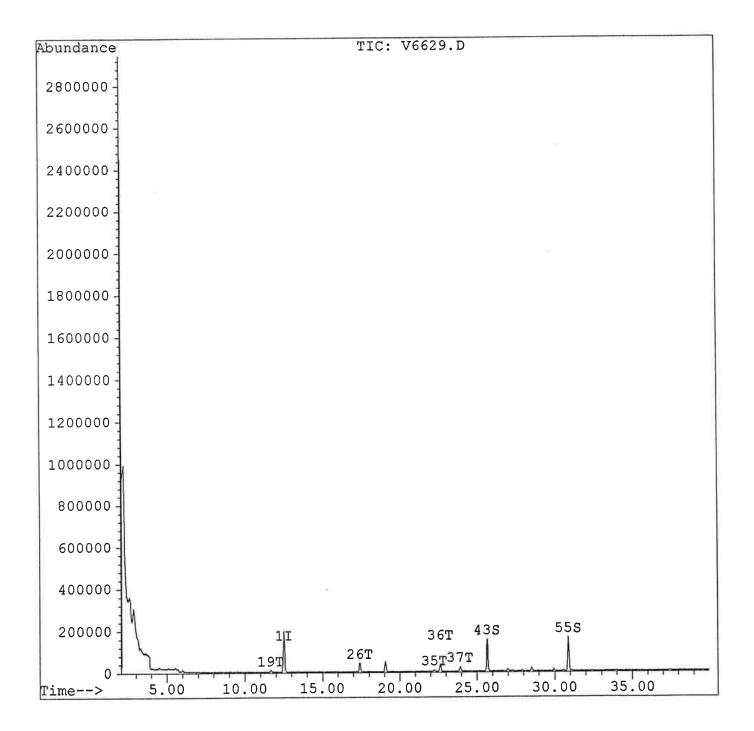
Misc : Lyondell - RC-EL-00-1099

Multiplr: 1.00

Quant Time: Oct 22 11:35 1999

Method : C:\HPCHEM\1\METHODS\RUN524.M

Title : 524.2 Purgable Organics
Last Update : Thu Oct 21 13:22:11 1999
Response via : Multiple Level Calibration



Data File : C:\HPCHEM\1\DATA\V6630.D Acq On : 21 Oct 99 5:27 pm

Vial: 14 Operator: vb

Sample : R-6204.5 Misc : Lyondell - RC-EC-69-1099

Inst : 5971 - In Multiplr: 1.00

Quant Time: Oct 22 11:33 1999

Method : C:\HPCHEM\1\METHODS\RUN524.M
Title : 524.2 Purgable Organics Last Update : Thu Oct 21 13:22:11 1999 Response via : Multiple Level Calibration

Internal Standards	R.T.	QIon	Response	Conc Units	Dev(Min)
1) Fluorobenzene	12.53	96	499715	5.00 ug/L	0.00
System Monitoring Compounds				8R	ecovery
43) 4-bromofluorobenzene	25,67	95	191795	5.11 ug/L	102.11%
55) 1,2-dichlorobenzene-d4	30.88	152	128748	5.32 ug/L	106.38%
Target Compounds					Qvalue
19) Benzene	11.68	78	36335	0.20 ug/L	96
26) Toluene	17.40	91	94149	0.60 ug/L	98
36) m&p-xylenes	22.64	106	35481	0.29 ug/L	92
37) o-xylene	23.94	91	45874	0.38 ug/L	93

<sup>(#) =</sup> qualifier out of range (m) = manual integration V6630.D RUN524.M Fri Oct 22 11:34:01 1999

Vial: 14

: 5971 - In

Operator: vb

Multiplr: 1.00

Inst

Data File : C:\HPCHEM\1\DATA\V6630.D

: 21 Oct 99 5:27 pm Acq On

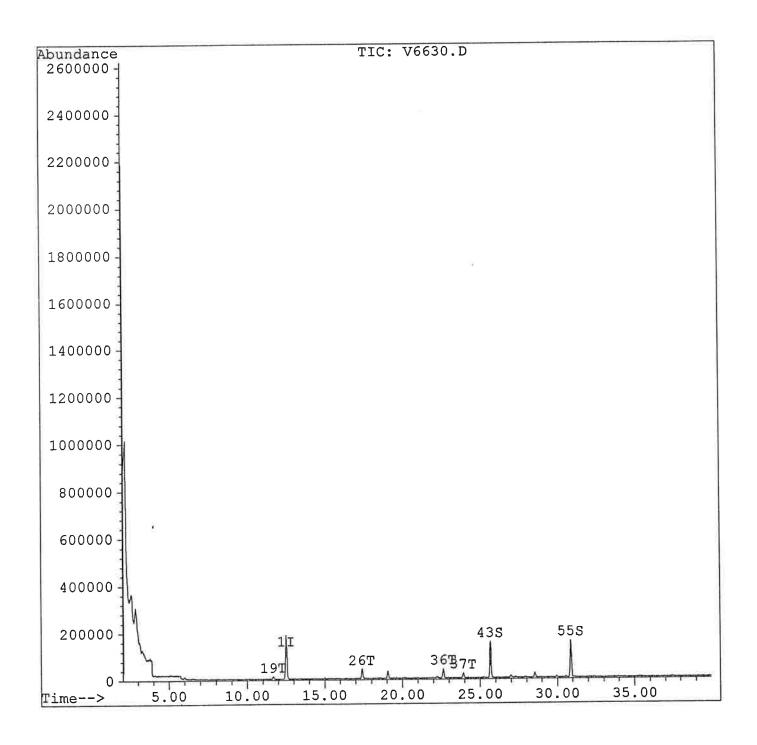
: R-6204.5 Sample

: Lyondell - RC-EC-69-1099

Misc Quant Time: Oct 22 11:33 1999

: C:\HPCHEM\1\METHODS\RUN524.M Method

: 524.2 Purgable Organics Title Last Update : Thu Oct 21 13:22:11 1999 Response via : Multiple Level Calibration



Vial: 15 Data File : C:\HPCHEM\1\DATA\V6631.D Acq On : 21 Oct 99 6:14 pm Operator: vb

Inst : 5971 - In Sample : R-6204.6 Misc : Lyondell - RC-EC-33-1099

Multiplr: 1.00

Quant Time: Oct 22 11:31 1999

: C:\HPCHEM\1\METHODS\RUN524.M Method

Title : 524.2 Purgable Organics
Last Update : Thu Oct 21 13:22:11 1999 Response via : Multiple Level Calibration

Internal Standards	R.T.	QIon	Response	Conc Units	Dev(Min)
1) Fluorobenzene	12.51	96	512975	5.00 ug/L	-0.02
System Monitoring Compounds 43) 4-bromofluorobenzene 55) 1,2-dichlorobenzene-d4	25.65 30.87	95 152	198194 131175	%! 5.14 ug/L 5.28 ug/L	
Target Compounds 19) Benzene 26) Toluene 35) Ethylbenzene 36) m&p-xylenes 37) o-xylene	11.67 17.40 22.22 22.63 23.91	78 91 91 106 91	79680 351516 77328 147430 168235	0.42 ug/L 2.18 ug/L 0.43 ug/L 1.19 ug/L 1.35 ug/L	Qvalue 94 96 95 97

<sup>(#) =</sup> qualifier out of range (m) = manual integration V6631.D RUN524.M Fri Oct 22 11:31:31 1999

Vial: 15 Operator: vb

Multiplr: 1.00

Inst : 5971 - In

Data File : C:\HPCHEM\1\DATA\V6631.D
Acg On : 21 Oct 99 6:14 pm

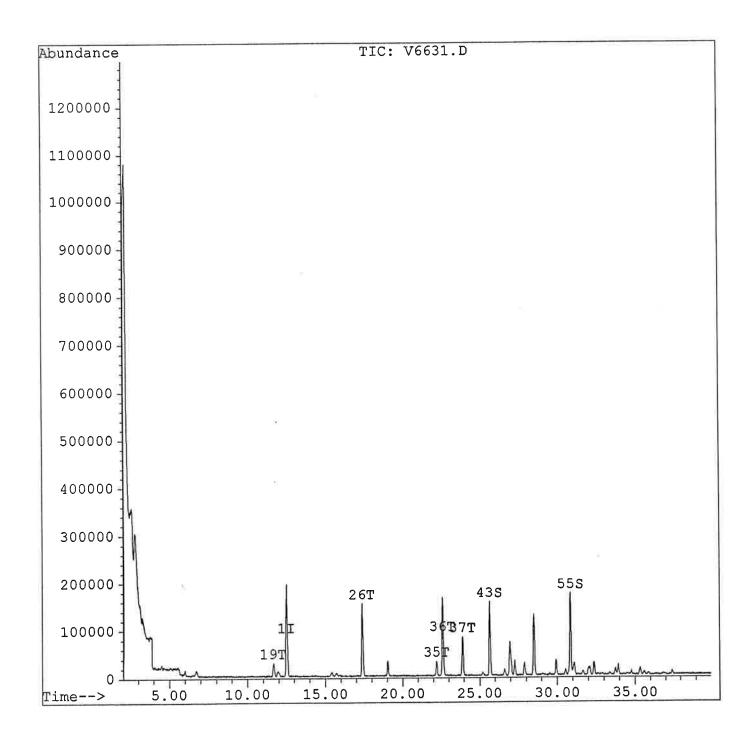
Sample : R-6204.6

Sample : R-6204.6 Misc : Lyondell - RC-EC-33-1099

Quant Time: Oct 22 11:31 1999

Method : C:\HPCHEM\1\METHODS\RUN524.M

Title : 524.2 Purgable Organics
Last Update : Thu Oct 21 13:22:11 1999
Response via : Multiple Level Calibration



### Quant. ation Report

Data File : C:\HPCHEM\1\DATA\V6632.D Vial: 16 Acq On : 21 Oct 99 Operator: vb 7:00 pm

Inst : 5971 - In Sample : R-6204.7

Multiplr: 1.00 Misc : Lyondell - RC-EC-00-1099 Quant Time: Oct 22 11:29 1999

: C:\HPCHEM\1\METHODS\RUN524.M Method

: 524.2 Purgable Organics Title Last Update : Thu Oct 21 13:22:11 1999 Response via : Multiple Level Calibration

Internal Standards	R.T.	QIon	Response	Conc Units	Dev(Min)
1) Fluorobenzene	12.52	96	530075	5.00 ug/L	-0.01
System Monitoring Compounds 43) 4-bromofluorobenzene 55) 1,2-dichlorobenzene-d4	25.67 30.87	95 152	198142 131666	%I 4.97 ug/L 5.13 ug/L	Recovery 99.44% 102.56%
Target Compounds 19) Benzene 26) Toluene 35) Ethylbenzene 36) m&p-xylenes 37) o-xylene	11.67 17.39 22.21 22.64 23.92	78 91 91 106 91	84522 414768 88610 186555 211986	0.43 ug/L 2.49 ug/L 0.48 ug/L 1.46 ug/L 1.65 ug/L	Qvalue 98 94 89 98

<sup>(#) =</sup> qualifier out of range (m) = manual integration V6632.D RUN524.M Fri Oct 22 11:29:35 1999

Vial: 16

Multiplr: 1.00

: 5971 - In

Operator: vb

Inst

Data File : C:\HPCHEM\1\DATA\V6632.D

Acq On : 21 Oct 99 7:00 pm

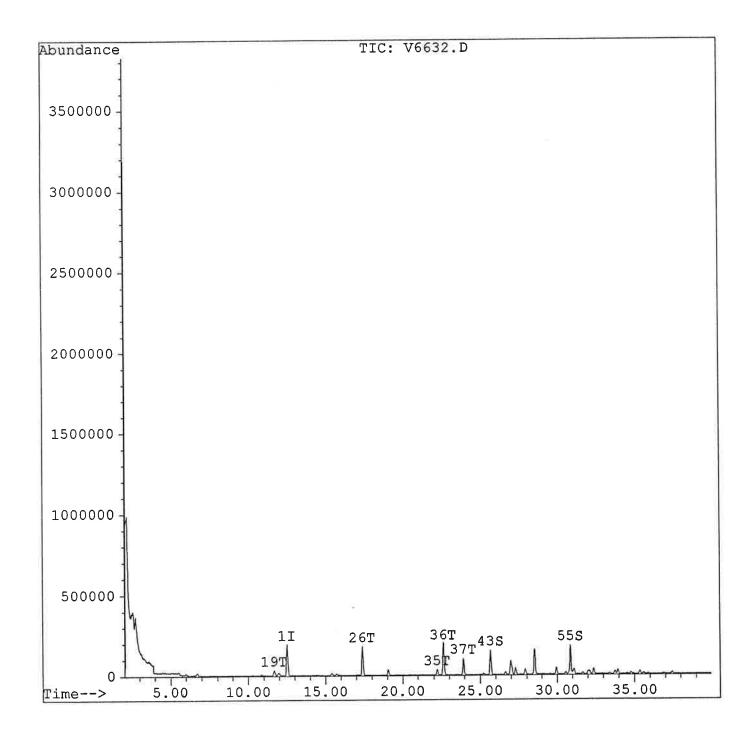
Sample : R-6204.7

Misc : Lyondell - RC-EC-00-1099

Quant Time: Oct 22 11:29 1999

Method : C:\HPCHEM\1\METHODS\RUN524.M

Title : 524.2 Purgable Organics
Last Update : Thu Oct 21 13:22:11 1999
Response via : Multiple Level Calibration



Data File : C:\HPCHEM\1\DATA\V6626.D Vial: 10 Operator: vb

Acq On : 21 Oct 99 2:21 pm Inst : 5971 - In : R-6204.8 Sample

Multiplr: 1.00 Misc : Lyondell - RC-EL-00-1099A

Quant Time: Oct 22 11:39 1999

: C:\HPCHEM\1\METHODS\RUN524.M Method

Title Last Update : Thu Oct 21 13:22:11 1999 Response via : Multiple Level Calibration

Internal Standards	R.T.	QIon	Response	Conc Units	Dev(Min)
1) Fluorobenzene	12.51	96	489543	5.00 ug/L	-0.02
System Monitoring Compounds 43) 4-bromofluorobenzene 55) 1,2-dichlorobenzene-d4	25.68 30.87	95 152	181678 120257	%I 4.94 ug/L 5.07 ug/L	Recovery 98.73% 101.43%
Target Compounds 19) Benzene 26) Toluene 35) Ethylbenzene 36) m&p-xylenes 37) o-xylene	11.67 17.41 22.23 22.65 23.92	78 91 91 106 91	38311 194136 43586 61715 75764	0.21 ug/L 1.26 ug/L 0.25 ug/L 0.52 ug/L 0.64 ug/L	Qvalue 97 98 98 91 97

<sup>(#) =</sup> qualifier out of range (m) = manual integration V6626.D RUN524.M Fri Oct 22 11:39:31 1999

Vial: 10

Inst : 5971 - In

Operator: vb

Multiplr: 1.00

Data File : C:\HPCHEM\1\DATA\V6626.D

Acq On : 21 Oct 99 2:21 pm

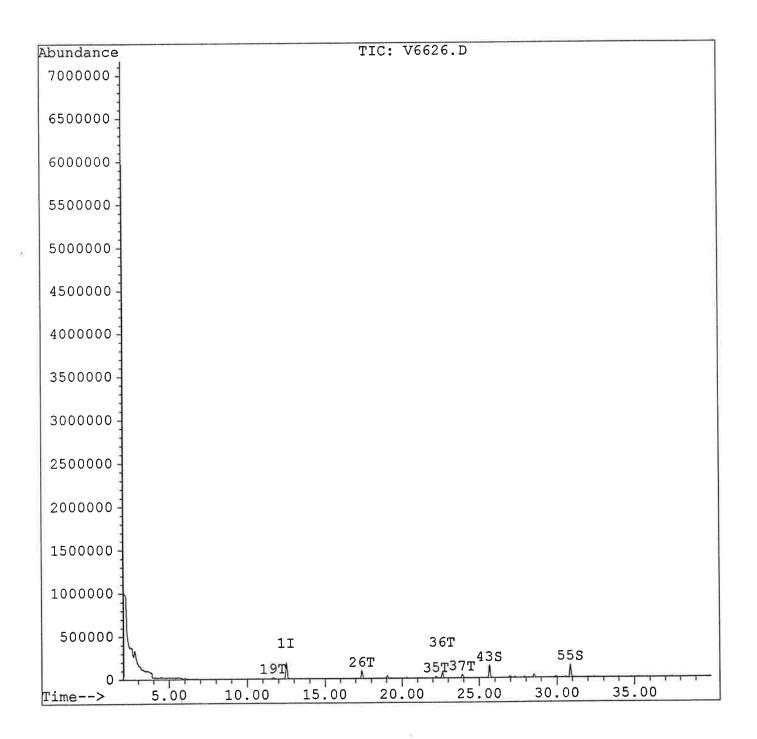
Sample : R-6204.8

Misc

: Lyondell - RC-EL-00-1099A

Quant Time: Oct 22 11:39 1999

Method : C:\HPCHEM\1\METHODS\RUN524.M



Data File : C:\HPCHEM\1\DATA\V6634.D

Acq On : 21 Oct 99 8:32 pm

Vial: 4 Operator: vb

Inst : 5971 - In Multiplr: 1.00

Sample : R-6204.9 Misc : Lyondell - Rinsate Blank Quant Time: Oct 22 9:10 1999

Method : C:\HPCHEM\1\METHODS\RUN524.M

Internal Standards	R.T.	QIon	Response	Conc Units	Dev(Min)
1) Fluorobenzene	12.50	96	496550	5.00 ug/L	-0.03
System Monitoring Compounds 43) 4-bromofluorobenzene 55) 1,2-dichlorobenzene-d4	25.64 30.85	95 152	186984 122328	%R 5.01 ug/L 5.09 ug/L	ecovery 100.18% 101.72%
Target Compounds 15) Chloroform 24) Bromodichloromethane 31) Dibromochloromethane	10.22 15.08 19.89	83 83 129	1336164 1277969 577469	15.93 ug/L 27.70 ug/L 28.53 ug/L	Qvalue 99 98 99

<sup>(#) =</sup> qualifier out of range (m) = manual integration V6634.D RUN524.M Fri Oct 22 11:26:11 1999

Vial: 4

Multiplr: 1.00

: 5971 - In

Operator: vb

Inst

Data File : C:\HPCHEM\1\DATA\V6634.D : 21 Oct 99 8:32 pm Acq On

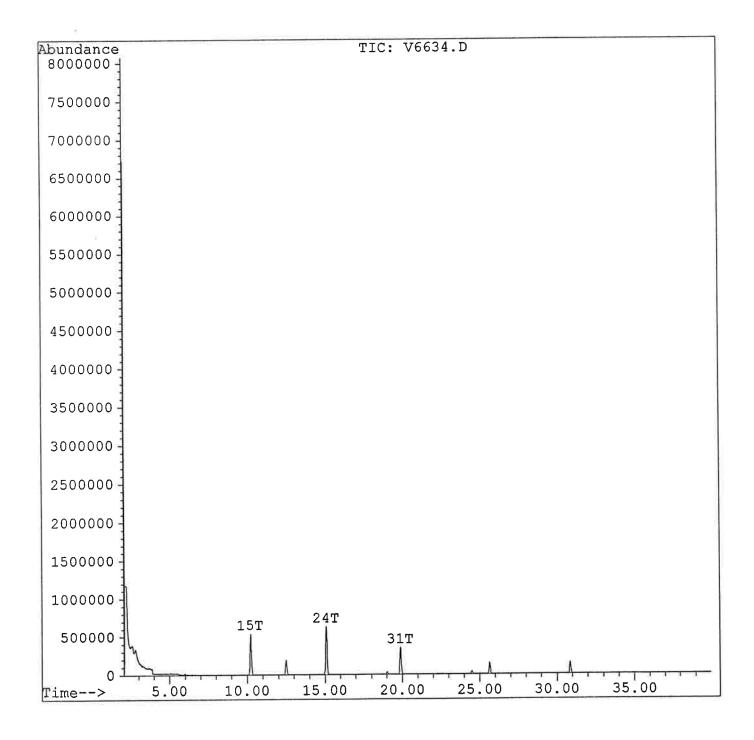
: R-6204.9 Sample

: Lyondell - Rinsate Blank

Misc

Quant Time: Oct 22 9:10 1999

: C:\HPCHEM\1\METHODS\RUN524.M Method



Data File : c:\hpchem\1\data\v6635.d

Acq On : 21 Oct 99 9:19 pm

Vial: 5 Operator: vb

Sample : R-6204.10 Misc : Lyondell - Trip Blank

Inst : 5971 - In

Multiplr: 1.00

Quant Time: Oct 22 9:10 1999

Last Update : Thu Oct 21 13:22:11 1999 Response via : Multiple Level Calibration

Internal Standards	R.T.	QIon	Response	Conc Units Dev(Min)
1) Fluorobenzene	12.50	96	449348	5.00 ug/L -0.03
System Monitoring Compounds 43) 4-bromofluorobenzene 55) 1,2-dichlorobenzene-d4	25.62 30.80	95 152	175205 113226	%Recovery 5.19 ug/L 103.73% 5.20 ug/L 104.04%
Target Compounds				Qvalue

Vial: 5
Operator: vb

Multiplr: 1.00

Inst : 5971 - In

Data File : c:\hpchem\1\data\v6635.d

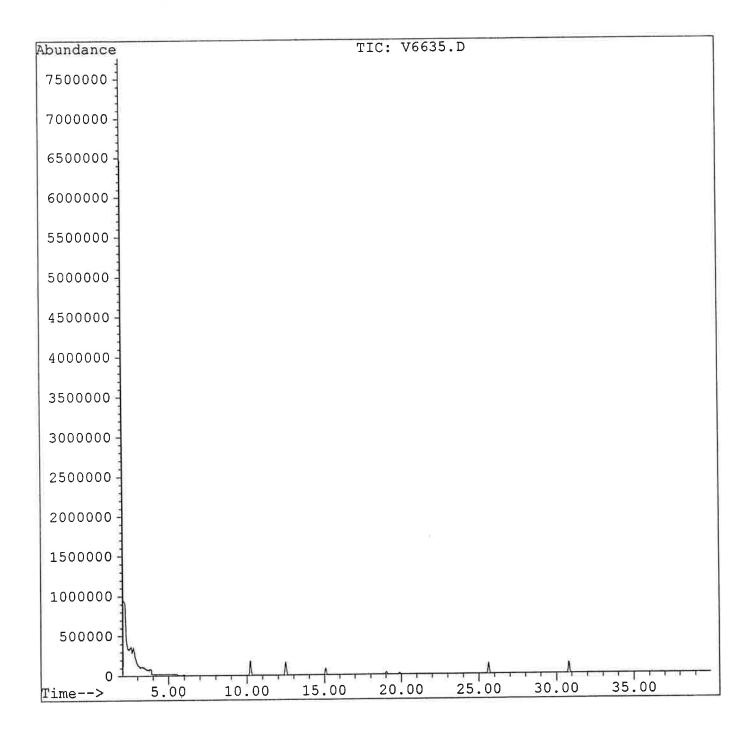
Acq On : 21 Oct 99 9:19 pm

Sample : R-6204.10

Misc : Lyondell - Trip Blank

Quant Time: Oct 22 9:10 1999

Method : C:\HPCHEM\1\METHODS\RUN524.M



# WATER VOLATILE SYSTEM MONITORING COMPOUND RECOVERY

Customer : Lyondel

Г		SMC1	SMC2		OTHER	TOT
	SAMPLE NO.	#	#	#	#	OUT
01	VBLK01	108				
02	R-6204.8	99				
03	R-6204.2	103				
04	R-6204.3	98				
05	R-6204.4	101				
06	R-6204.5	102				$\vdash$
07	R-6204.6	103				
08	R-6204.7	99				
09	R-6204.1	100				_
10	R-6204.9	100				-
11	R-6204.10	104			<u> </u>	-
12	VBLK02	93				-
13	R-6204.6MS	96				
14	R-6204.6MSD	99			1	
15						-
16					-	-
17						
18						+
19						+
20						_
21						+
22					-	
23						-
24					-	+
25						-
26				-	_	-
27						+
28				-		
29				1		-
30						

SMC1 = 4-Bromofluorobenzene

SMC2 = 1,2-dichlorobenzene-d4

QC LIMITS (75-115) (75-115)

- # Column to be used to flag recovery values
- Values outside of contract required QC limits
- D System Monitoring Compound diluted out

FORM II VOA-1

# RELIANCE LABORATORIES, INC. WATER SEMIVOLATILE MATRIX SPIKE/MATRIX SPIKE DUPLICATE RECOVERY

Matrix Spike - Sample No.: R-6204.6

ASHAN ASHAN	SPIKE ADDED	SAMPLE CONCENTRATION	MS CONCENTRATION	MS %	QC. LIMITS
COMPOUND	(ug/Kg)	(ug/Kg)	(ug/Kg)	REC #	REC.
Benzene	3.00	0.21	3.27	102	(80-120)
Toluene	3.00	1.09	4.05	99	(80-120)
Ethylbenzene	3.00	0.22	3.32	103	(80-120)
m&p-xylenes	3.00	0.59	3.77	106	(80-120)
o-xylenes	3.00	0.68	3.80	104	(80-120)
Styrene	3.00	0.00	3.06	102	(80-120)

	SPIKE	MSD	MSD			
	ADDED	CONCENTRATION	%	%	QC L	IMITS
COMPOUND	(ug/Kg)	(ug/Kg)	REC #	RPD #	RPD	REC.
Benzene	3.00	3.15	98	4	20	(80-120)
Toluene	3.00	3.96	97	2	20	(80-120)
Ethylbenzene	3.00	3.18	99	5	20	(80-120)
m&p-xylenes	3.00	3.58	100	6	20	(80-120)
o-xylenes	3.00	3.69	100	4	20	(80-120)
Styrene	3.00	2.83	94	8	20	(80-120)

# Column to be used to flag recovery and RPD values with an asteris														
# Calumn to be used to tiga recovery and PPH Values with an asiens	k	asteris	an	with	values	PPD	hae	racovani	floor	ucod to	ho	mn to	Calu	ш

Comments: _	

<sup>#</sup> Column to be used to flag re\* Values outside of QC limits

#### VOLATILE METHOD BLANK SUMMARY

VBLK0

	Customer :	Lyondel			
Lab File ID V66	25.D			Lab Sample ID:	BLANK1
Date Analyzed:	10/21/99			Time Analyzed:	1334
GC Column:	DB-624 ID:	0.53(mm)			
Instrument ID:	HP5971				
	S METHOD BLANK	APPLIES TO THE	FOLLOWING :	SAMPLES, MS A	ND MSD:
		LAB	LAB	TIME	
	SAMPLE NO.	SAMPLE ID	FILE ID	ANALYZED	
01	R-6204.8	EL00A	V6626.D	1421	
	R-6204.2	ER00	V6627.D	1508	
03	R-6204.3	EL19	V6628.D	1554	
04	R-6204.4	EL00	V6629.D	1641	
05	R-6204.5	EC69	V6630.D	1727	
06	R-6204.6	EC33	V6631.D	1814	
	R-6204.7	EC00	V6632.D	1900	
	R-6204.1	ER65	V6633.D	1946	
09	R-6204.9	RBLK	V6634.D	2032	
	R-6204.10	TBLK	V6635.D	2119	
11					
12					
13					
14					
15					
16					
17					
18					
19					
20					
21					
22					
23					

COMMENTS:			
			_

Page 1 of 1

FORM IV VOA

Data File : C:\HPCHEM\1\DATA\V6625.D

Vial: 9

Acq On : 21 Oct 99 1:34 pm

Operator: vb

Inst : 5971 - In

Multiplr: 1.00

Sample : VBLK01
Misc : Method Blank Quant Time: Oct 22 11:40 1999

Method : C:\HPCHEM\1\METHODS\RUN524.M
Title : 524.2 Purgable Organics

Last Update : Thu Oct 21 13:22:11 1999

Response via : Multiple Level Calibration

Internal Standards	R.T.	QIon	Response	Conc Units	Dev(Min)
1) Fluorobenzene	12.52	96	515565	5.00 ug/L	0.00
System Monitoring Compounds 43) 4-bromofluorobenzene 55) 1,2-dichlorobenzene-d4	25.67 30.87	95 152	209884 141082	%R 5.42 ug/L 5.65 ug/L	
Target Compounds					Qvalue

<sup>(#) =</sup> qualifier out of range (m) = manual integration V6625.D RUN524.M Fri Oct 22 11:40:15 1999

Vial: 9

Multiplr: 1.00

: 5971 - In

Operator: vb

Inst

Data File : C:\HPCHEM\1\DATA\V6625.D

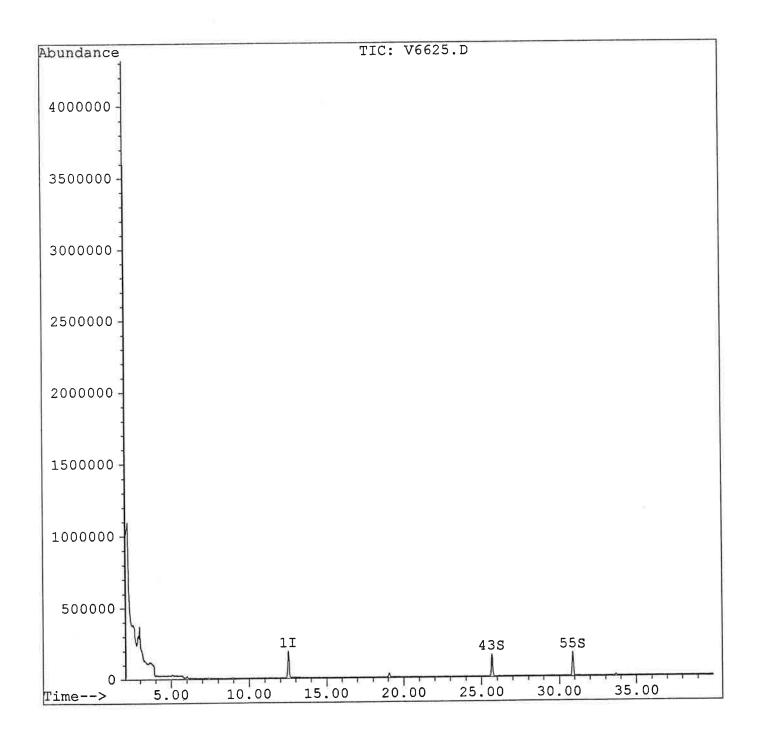
Acq On : 21 Oct 99 1:34 pm

Sample : VBLK01

Misc : Method Blank

Quant Time: Oct 22 11:40 1999

Method : C:\HPCHEM\1\METHODS\RUN524.M



#### VOLATILE METHOD BLANK SUMMARY

VBLK0

Lab File ID <u>V6638</u>				Lab Sample ID	•
Date Analyzed: 1	0/22/99			Time Analyzed	. 1002
GC Column: D	B-624 ID	: <u>0.53</u> (mm	)		
Instrument ID: H	P5971				
THIS	METHOD BLANK	APPLIES TO THE	FOLLOWING	SAMPLES, MS A	ND MSD:
ir-		LAB	LAB	TIME	
	SAMPLE NO.	SAMPLE ID	FILE ID	ANALYZED	
01 R	-6204.6MS	EC33MS	V6639.D	1650	
	-6204.6MSD	EC33MSD	V6640.D	1737	
03			ļ		
04			<u> </u>	1	
05 06					
07				<u> </u>	
08					
09					
10					
11					
12			-	-	
13					
14 15					•
16					1
17					]
18					]
19					
20					ł
21					1
22 23			<del> </del>	-	1
23					1
25					
26					]
27					1
28					-
29		<del></del>		-	1
30					1
COMMENTS:					

FORM IV VOA

Vial: 9

Multiplr: 1.00

: 5971 - In

Operator: vb

Inst

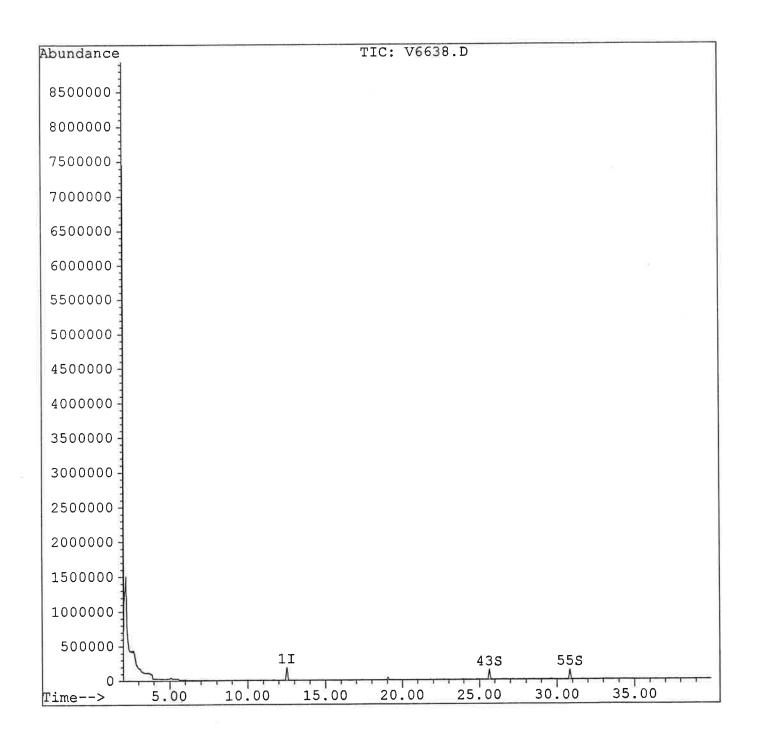
Data File : C:\HPCHEM\1\DATA\V6638.D

Acq On : 22 Oct 99 4:02 pm

Sample : Method Blank
Misc : Method Blank

Quant Time: Oct 27 9:48 1999

Method : C:\HPCHEM\1\METHODS\RUN524.M



#### VOLATILE ORGANIC INSTRUMENT PERFORMANCE CHECK BROMOFLUOROBENZENE (BFB)

Customer : Lyondel	
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Lab File ID: V6621.D

BFB Injection Date: 10/21/99

Instrument ID: HP5971A

BFB Injection Time: 0945

GC Column: DB-624

ID: 0.53 (mm)

	%RELATIVE
ION ABUNDANCE CRITERIA	ABUNDANCE
8.0 - 40.0% of mass 95	22.2
30.0 - 66.0% of mass 95	47.4
Base peak, 100% relative abundance	100.0
5.0 - 9.0% of mass 95	6.7
Less than 2.0% of mass 174	0.0 ( 0.0 )1
50.0 - 120.0% of mass 95	65.1
4.0 - 9.0% of mass 174	5.2 ( 8.0 )1
93.0 - 101.0% of mass 174	62.4 ( 96.0 )1
	3.8 ( 6.1 )2
	8.0 - 40.0% of mass 95 30.0 - 66.0% of mass 95 Base peak, 100% relative abundance 5.0 - 9.0% of mass 95 Less than 2.0% of mass 174 50.0 - 120.0% of mass 95 4.0 - 9.0% of mass 174

1-Value is % mass 174

2-Value is % mass 176

### This check applies to the following SAMPLES, MS, MSD, BLANKS and STANDARDS:

		LAB	LAB	DATE	TIME
	SAMPLE NO.	SAMPLE ID	FILE ID	ANALYZED	ANALYZED
01	VSTD050	ICC005	V6622.D	10/21/99	1033
02	VSTD020	ICC002	V6623.D	10/21/99	1120
03	VSTD010	ICC001	V6624.D	10/21/99	1217
04	VBLK01	BLANK1	V6625.D	10/21/99	1334
05	R-6204.8	EL00A	V6626.D	10/21/99	1421
	R-6204.2	ER00	V6627.D	10/21/99	1508
07	R-6204.3	EL19	V6628.D	10/21/99	1554
	R-6204.4	EL00	V6629.D	10/21/99	1641
	R-6204.5	EC69	V6630.D	10/21/99	1727
	R-6204.6	EC33	V6631.D	10/21/99	1814
	R-6204.7	EC00	V6632.D	10/21/99	1900
	R-6204.1	ER65	V6633.D	10/21/99	1946
	R-6204.9	RBLK	V6634.D	10/21/99	2032
	R-6204.10	TBLK	V6635.D	10/21/99	2119
15					
16					
17					
18					
19					
20					
21					
22					

Page 1 of 1

FORM V VOA

# VOLATILE ORGANIC INSTRUMENT PERFORMANCE CHECK BROMOFLUOROBENZENE (BFB)

Customer: Lyondei	
Lab File ID: V6636.D	BFB Injection Date: 10/22/99
Instrument ID: HP5971A	BFB Injection Time: 1438

GC Column: DB-624 ID: 0.53 (mm)

		%RELATIVE
m/e	ION ABUNDANCE CRITERIA	ABUNDANCE
50	8.0 - 40.0% of mass 95	22.0
75	30.0 - 66.0% of mass 95	45.1
95	Base peak, 100% relative abundance	100.0
96	5.0 - 9.0% of mass 95	6.5
173	Less than 2.0% of mass 174	0.0 ( 0.0 )1
174	50.0 - 120.0% of mass 95	63.7
175	4.0 - 9.0% of mass 174	4.7 ( 7.4 )1
176	93.0 - 101.0% of mass 174	63.8 ( 100.3 )1
177	5.0 - 9.0% of mass 176	4.4 ( 6.8)2

1-Value is % mass 174 2-Value is % mass 176

This check applies to the following SAMPLES, MS, MSD, BLANKS and STANDARDS:

		LAB	LAB	DATE	TIME
	SAMPLE NO.	SAMPLE ID	FILE ID	ANALYZED	ANALYZED
01	VSTD050	CC002	V6637.D	10/22/99	1510
02	VBLK02	BLANK2	V6638.D	10/22/99	1602
03	R-6204.6MS	EC33MS	V6639.D	10/22/99	1650
04	R-6204.6MSD	EC33MSD	V6640.D	10/22/99	1737
05					
06					
07					
80					
09					
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19					<b>.</b>
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21					
22				1	

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FORM V VOA

Data File : C:\HPCHEM\1\DATA\V6636.D

Acq On : 22 Oct 99 2:38 pm

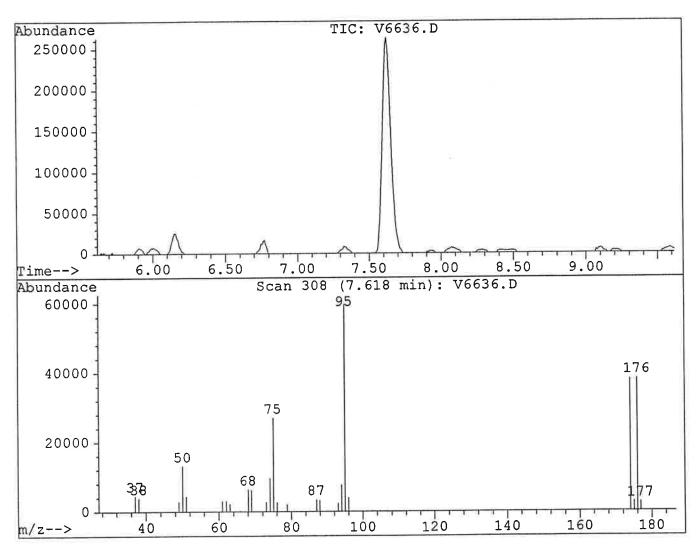
Sample : bfb Misc : bfb Vial: 1
Operator: vb

Inst : 5971 - In

Multiplr: 1.00

Method : C:\HPCHEM\1\METHODS\RUN524.M

Title : 524.2 Purgable Organics



Peak Apex is scan: 308

1	Target Mass		Rel. to Mass	1	Lower Limit%	1	Upper Limit%		Rel. Abn%		Raw Abn		Result Pass/Fail	
1	50	 I	 95	1	15	Ĭ	40	1	22.0	1	13105	١	PASS	1
i	75	i	95	ì	30	Î	80	Î	45.1	1	26944	1	PASS	1
i	95	i	95	î	100	ì	100	1	100.0	- 1	59680	1	PASS	1
ì	96	i	95	î	5	i	9	Ť	6.5	1	3901	1	PASS	1
÷	173	î	174	i	0	i	2	Ì	0.0	1	0	1	PASS	1
i	174	ıi.	95	î	50	i	100	ì	63.7	1	37992	1	PASS	1
ï	175	i	174	Î	5	Î	9	î	7.4	1	2805	1	PASS	1
i	176	i	174	i	95	i	101	Î	100.3	Ì	38096	1	PASS	1
i	177	i	176	i	5	j	9	i	6.8	Ĺ	2607	1	PASS	1

#### **VOLATILE ORGANICS INITIAL CALIBRATION DATA**

Customer	Lyondel	
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Instrument ID: <u>HP5971A</u> Calibration Date(s): <u>10/21/99</u> <u>10/21/99</u>

Calibration Times: 1033 1217

GC Column: DB-624 ID: <u>0.53</u> (mm)

Lab File ID:	RRF01 =	V6624.D	R	RF02 = V6	623.D	
RRF05 = V6622.D						%
COMPOUND	RRF01	RRF02	RRF05		RRF	RSE
Benzen <b>e</b>	1.803	1.783	1.923		1.836	4.1
Toluene	1.443	1.620	1.646		1.570	7.0
Ethylbenzene	1.534	1.869	1.859		1.754	10.9
m&p-xylenes	1.039	1.294	1.285		1.206	12.0
o-xylene	1.073	1.269	1.299		1.214	10.1
Styrene	0.746	0.891	0.930		0.856	11.3
4-bromofluorobenzene	0.371	0.387	0.370		0.376	2.5
1,2-dichlorobenzene-d4	0.241	0.241	0.245		0.242	1.0

Page 1 of 1

FORM VI VOA

#### VOLATILE CONTINUING CALIBRATION CHECK

Customer : Lyondell

Instrument ID: HP5971A Calibration Date: 10/22/99

Time: 1225

Lab File ID: V6298.D Init. Calib. Date(s): 10/21/99 10/21/99

Init. Calib. Times: 1033 1217

GC Column:

DB-624

ID: <u>0.53</u> (mm)

			MIN		MAX
COMPOUND	RRF	RRF20	RRF	%D	%D
Benzene	1.837	1.875		-2.1	
Toluene	1.570	1.647		-4.9	
Ethylbenzene	1.754	1.834		-4.6	
m&p-xylenes	1.206	1.266		-5.0	
o-xylene	1.213	1.279		-5.4	
Styrene	0.856	0.883		-3.2	
					-
					-
		-			-
					-
4-Bromofluorobenzene	0.376	0.390		-3.7	
1,2-dichlorobenzene-d4	0.242	0.257		3.5	

All other compounds must meet a minimum RRF of 0.010.

Page 1 of 1

FORM VII VOA

#### VOLATILE INTERNAL STANDARD AREA AND RT SUMMARY

Customer: Lyondel

Lab File ID (Standard): V6622.D

Date Analyzed: 10/21/99

Instrument ID: HP5971A

Time Analyzed: 1033

GC Column: DB-624

ID: 0.53 (mm)

	IS1					T				
	AREA #	RT #	AREA	#	RT	#	AREA	#	RT	#
12 HOUR STD	561288	12.53								
UPPER LIMIT	1122576	13.03								
LOWER LIMIT	280644	12.03								
SAMPLE										
NO.										
VBLK01	515565	12.52								
)2 R-6204.8	489543	12.51								-
3 R-6204.2	487047	12.52								
04 R-6204.3	473211	12.52								
05 R-6204.4	524441	12.52								
06 R-6204.5	499715	12.53								
07 R-6204.6	512975	12.51								_
08 R-6204.7	530075	12.52								_
09 R-6204.1	506914	12.49								
10 R-6204.9	496550	12.50								_
11 R-6204.10	449348	12.50								
12										
13									<u> </u>	
14									-	
15								_	ļ	_
16										
17										
18									-	_
19										_
20										
21										
22										

IS1 = Fluorobenzene

AREA UPPER LIMIT = +100% of internal standard area AREA LOWER LIMIT = -50% of internal standard area RT UPPER LIMIT = +0.50 minutes of internal standard RT RT LOWER LIMIT = -0.50 minutes of internal standard RT

# Column used to flag values outside QC limits with an asterisk

\* Values outside of QC limits.

Page 1 of 1

FORM VIII VOA

# CHAIN OF CUSTODY

Hydrology Applied Associates, Inc.

Date:

Page 1 of 1

LAB ID: R-6204

Trun around time Standard ... (stanrard / rush)

Project ID:

400 Frankfort Rd, Monaca, Pa 15061

Sampler Name: BDP
Preserved Y 🖎
Sample Intact: 🕎 N

Address:

724) 728 - 6586

724) 728 - 6498

Customer:

Lyondell Chemical

00 Frankfort Rd Monaca Pa. 15061

E-Mail: Fax: Phone

b.petroff@worldnet.att.net

# of containers Water MATRIX Soil Other EPA Test Method 524.2 BTEX (602/8020) TPH (418.1) VOA (624/8260) + 15 ORGANICS BNA / BN / + 15 Pest / Herb PCB's TCLP Organics / pp + 40 Other TCLP / RCRA (8) Priority Pol. (13) METALS Total Metals (list below) Dissolved Metals Other pH / CN / Sulfide Fl pt / % Solids OTHERS O & G / TSS / TOX BOD / COD / TOC E-Mail results (Y) N Fax results: Q N

RC-ER-65-1099 Sample ID RC-EL-00-1099A RC-EC-69-1099 RC+EL-00-1099 RC-EL-19-1099 RC-ER-00-1099 RC-EC-00-1099 RC-EC-33-1099 Rinsate Blank nstructions: Please Fax results to and Skip Meier: (303) 873 - 6110 and E-Mail results to Brian Petroff: b.petroff@worldnet.att.net. rip Blank Date Spld. 10/20/99 10/20/99 10/20/99 10/20/99 10/20/99 10/20/99 10/20/99 10/20/99 10/20/99 10/20/99 Time Spld 10:52 10:47 11:03 11:35 10:52 10:58 10:57 10:56 11:02 9:45 1 × × × ××

Agent of: AHA
Received by:
Rece

Date / Time: 10/2/64 740

Date / Time: Agent of: Received by: Agent of:

Agent of:

Date / Time:

Agent of: Received by:

Submitted by:\_

Submitted by: BDP

Submitted by:\_

			Deliverables:	Report to:
Customized	Reduced	Standard		ŀ

10/20/993:40 PM